

Connecting via winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal653adk

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *
* *

| | |
|----------------|---|
| NEWS 1 | Web Page URLs for STN Seminar Schedule - N. America |
| NEWS 2 | "Ask CAS" for self-help around the clock |
| NEWS 3 SEP 01 | New pricing for the Save Answers for SciFinder |
| Wizard within | |
| | STN Express with Discover! |
| NEWS 4 OCT 28 | KOREAPAT now available on STN |
| NEWS 5 NOV 30 | PHAR reloaded with additional data |
| NEWS 6 DEC 01 | LISA now available on STN |
| NEWS 7 DEC 09 | 12 databases to be removed from STN on December 31, 2004 |
| NEWS 8 DEC 15 | MEDLINE update schedule for December 2004 |
| NEWS 9 DEC 17 | ELCOM reloaded; updating to resume; current-awareness |
| | alerts (SDIs) affected |
| NEWS 10 DEC 17 | COMPUAB reloaded; updating to resume; current-awareness |
| | alerts (SDIs) affected |
| NEWS 11 DEC 17 | SOLIDSTATE reloaded; updating to resume; current-awareness |
| | alerts (SDIs) affected |
| NEWS 12 DEC 17 | CERAB reloaded; updating to resume; current-awareness |
| | alerts (SDIs) affected |
| NEWS 13 DEC 17 | THREE NEW FIELDS ADDED TO IFIPAT/IFIUDB/IFICDB |
| NEWS EXPRESS | OCTOBER 29 CURRENT WINDOWS VERSION IS V7.01A, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004 |
| NEWS HOURS | STN Operating Hours Plus Help Desk Availability |
| NEWS INTER | General Internet Information |
| NEWS LOGIN | Welcome Banner and News Items |
| NEWS PHONE | Direct Dial and Telecommunication Network Access to STN |
| NEWS WWW | CAS world wide web site (general information) |

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer

agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 10:14:29 ON 28 DEC 2004

| | |
|----------------------|------------|
| => fil reg | |
| COST IN U.S. DOLLARS | SINCE FILE |
| TOTAL | ENTRY |
| SESSION | |
| FULL ESTIMATED COST | 0.21 |
| 0.21 | |

FILE 'REGISTRY' ENTERED AT 10:14:35 ON 28 DEC 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 26 DEC 2004 HIGHEST RN 802853-20-9
DICTIONARY FILE UPDATES: 26 DEC 2004 HIGHEST RN 802853-20-9

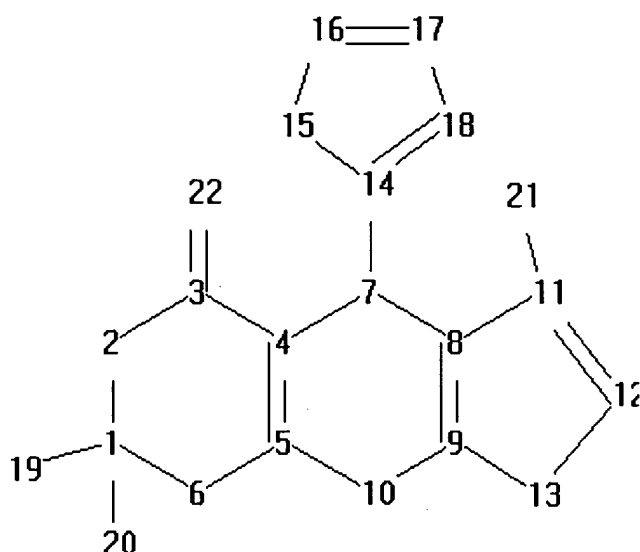
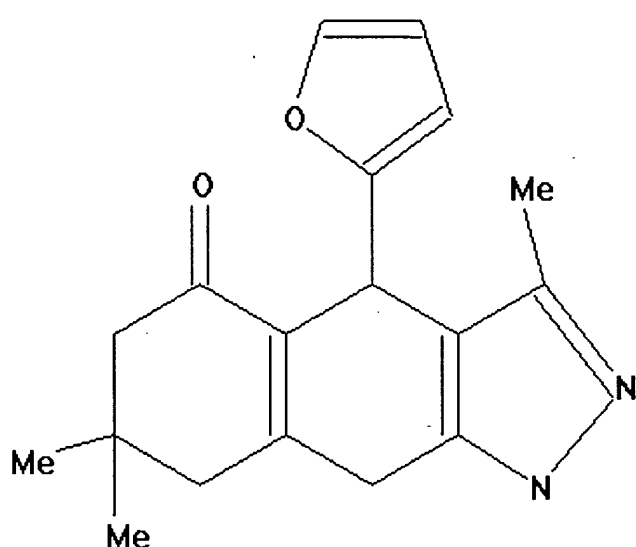
TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>
Uploading H:\STN queries\10612885.str



```

chain nodes :
19 20 21 22
ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18
chain bonds :
1-19 1-20 3-22 7-14 11-21
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10 9-13
11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact/norm bonds :
1-2 1-6 2-3 3-4 3-22 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10
9-13 11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact bonds :
1-19 1-20 7-14 11-21

```

```

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom
9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom
17:Atom 18:Atom 19:CLASS 20:CLASS 21:CLASS 22:CLASS

```

L1 STRUCTURE UPLOADED

=> s l1

SAMPLE SEARCH INITIATED 10:14:50 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 8 TO ITERATE

100.0% PROCESSED 8 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 8 TO 329
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s l1 ful

FULL SEARCH INITIATED 10:14:53 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 166 TO ITERATE

100.0% PROCESSED 166 ITERATIONS

0

ANSWERS

SEARCH TIME: 00.00.01

L3 0 SEA SSS FUL L1

=> s l1 ful fam

FULL SEARCH INITIATED 10:14:58 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS

0

ANSWERS

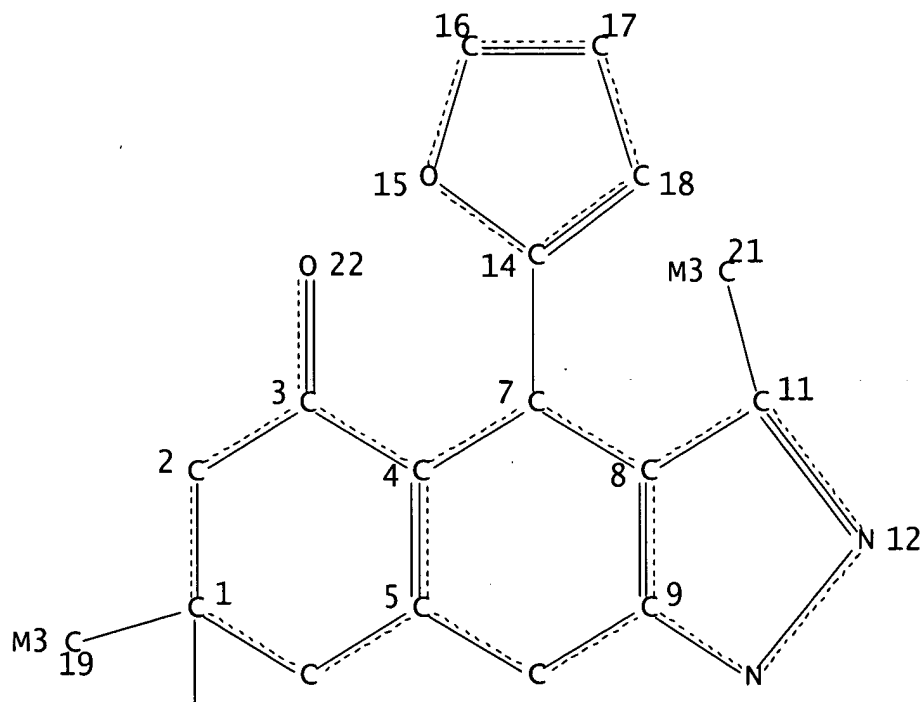
SEARCH TIME: 00.00.01

L4 0 SEA FAM FUL L1

=> d l1

L1 HAS NO ANSWERS

L1 STR



Page 2-A

NODE ATTRIBUTES:

| | | | | |
|--------|----|----|----|----|
| HCOUNT | IS | M3 | AT | 19 |
| HCOUNT | IS | M3 | AT | 20 |
| HCOUNT | IS | M3 | AT | 21 |
| NSPEC | IS | R | AT | 1 |
| NSPEC | IS | R | AT | 2 |
| NSPEC | IS | R | AT | 3 |
| NSPEC | IS | R | AT | 4 |
| NSPEC | IS | R | AT | 5 |
| NSPEC | IS | R | AT | 6 |
| NSPEC | IS | R | AT | 7 |
| NSPEC | IS | R | AT | 8 |
| NSPEC | IS | R | AT | 9 |
| NSPEC | IS | R | AT | 10 |
| NSPEC | IS | R | AT | 11 |
| NSPEC | IS | R | AT | 12 |
| NSPEC | IS | R | AT | 13 |
| NSPEC | IS | R | AT | 14 |
| NSPEC | IS | R | AT | 15 |
| NSPEC | IS | R | AT | 16 |
| NSPEC | IS | R | AT | 17 |
| NSPEC | IS | R | AT | 18 |
| NSPEC | IS | C | AT | 19 |
| NSPEC | IS | C | AT | 20 |
| NSPEC | IS | C | AT | 21 |
| NSPEC | IS | C | AT | 22 |

DEFAULT MLEVEL IS ATOM
MLEVEL IS CLASS AT 19 20 21 22
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

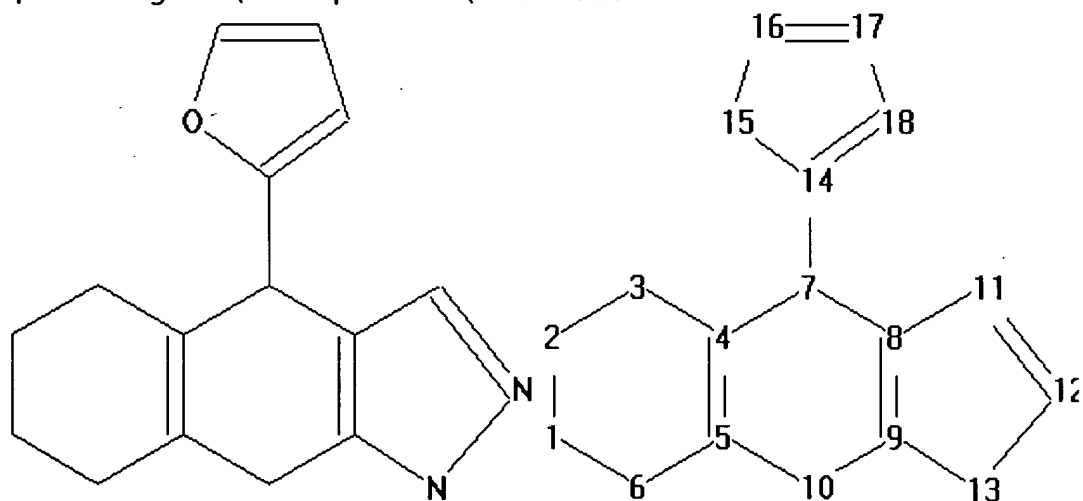
RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 22

STEREO ATTRIBUTES: NONE

=>

Uploading H:\STN queries\10612885a.str



```

ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18
chain bonds :
7-14
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10 9-13
11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact/norm bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10 9-13
11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact bonds :
7-14

```

```
Match level :
1:Atom  2:Atom  3:Atom  4:Atom  5:Atom  6:Atom  7:Atom  8:Atom
9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom
17:Atom 18:Atom
```

L5 STRUCTURE UPLOADED

```
=> s 15
SAMPLE SEARCH INITIATED 10:18:34 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -      24 TO ITERATE
```

```
100.0% PROCESSED      24 ITERATIONS      0
ANSWERS
SEARCH TIME: 00.00.01
```

```

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH    **COMPLETE**
PROJECTED ITERATIONS:   187 TO      773
PROJECTED ANSWERS:      0 TO      0

```

L6 0 SEA SSS SAM L5

```
=> s 15 fu1
FULL SEARCH INITIATED 10:18:38 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -      607 TO ITERATE
```

```
100.0% PROCESSED      607 ITERATIONS      0
ANSWERS
SEARCH TIME: 00.00.01
```

L7 0 SEA SSS FUL L5

```
=> s 15 fam
SAMPLE SEARCH INITIATED 10:18:42 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE .
```

```
100.0% PROCESSED      0 ITERATIONS      0
ANSWERS
SEARCH TIME: 00.00.01
```

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 0 TO 0
PROJECTED ANSWERS: 0 TO 0

L8 0 SEA FAM SAM L5

=> s 15 fam ful
FULL SEARCH INITIATED 10:18:48 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 2 TO ITERATE

100.0% PROCESSED 2 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

L9 0 SEA FAM FUL L5

=> DIS HIST

(FILE 'HOME' ENTERED AT 10:14:29 ON 28 DEC 2004)

FILE 'REGISTRY' ENTERED AT 10:14:35 ON 28 DEC 2004

L1 STRUCTURE UPLOADED
L2 0 S L1
L3 0 S L1 FUL
L4 0 S L1 FUL FAM
L5 STRUCTURE UPLOADED
L6 0 S L5
L7 0 S L5 FUL
L8 0 S L5 FAM
L9 0 S L5 FAM FUL

=>

---Logging off of STN---

=>
Executing the logoff script...

=> LOG Y

| COST IN U.S. DOLLARS | SINCE FILE |
|----------------------|------------|
| TOTAL | ENTRY |
| SESSION | |
| FULL ESTIMATED COST | 434.86 |
| 435.07 | |

STN INTERNATIONAL LOGOFF AT 10:19:23 ON 28 DEC 2004

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:sssptal653adk

PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

***** welcome to STN International *****

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 "Ask CAS" for self-help around the clock
NEWS 3 SEP 01 New pricing for the Save Answers for SciFinder
Wizard within
NEWS 4 OCT 28 STN Express with Discover!
NEWS 5 NOV 30 KOREAPAT now available on STN
NEWS 6 DEC 01 PHAR reloaded with additional data
NEWS 7 DEC 09 LISA now available on STN
2004 12 databases to be removed from STN on December 31,
NEWS 8 DEC 15 MEDLINE update schedule for December 2004
NEWS 9 DEC 17 ELCOM reloaded; updating to resume; current-
awareness
NEWS 10 DEC 17 alerts (SDIs) affected
awareness COMPUAB reloaded; updating to resume; current-
NEWS 11 DEC 17 alerts (SDIs) affected
awareness SOLIDSTATE reloaded; updating to resume; current-
NEWS 12 DEC 17 alerts (SDIs) affected
awareness CERAB reloaded; updating to resume; current-
NEWS 13 DEC 17 alerts (SDIs) affected
THREE NEW FIELDS ADDED TO IFIPAT/IFIUDB/IFICDB
NEWS EXPRESS OCTOBER 29 CURRENT WINDOWS VERSION IS V7.01A, CURRENT
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0j(c(JP),
AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004
STN Operating Hours Plus Help Desk Availability
NEWS HOURS General Internet Information
NEWS INTER General Internet Information
NEWS LOGIN Welcome Banner and News Items
NEWS PHONE Direct Dial and Telecommunication Network Access to
STN
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that
specific topic.

All use of STN is subject to the provisions of the STN Customer

agreement. Please note that this agreement limits use to
scientific research. Use for software development or design or implementation
of commercial gateways or other similar uses is prohibited and may
result in loss of user privileges and other penalties.

***** STN Columbus *****

FILE 'HOME' ENTERED AT 10:25:10 ON 28 DEC 2004

=> FILE REGISTRY
COST IN U.S. DOLLARS
TOTAL
ENTRY
SESSION
FULL ESTIMATED COST 0.21

FILE 'REGISTRY' ENTERED AT 10:25:22 ON 28 DEC 2004
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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provided by InfoChem.

STRUCTURE FILE UPDATES: 26 DEC 2004 HIGHEST RN 802853-20-9
DICTIONARY FILE UPDATES: 26 DEC 2004 HIGHEST RN 802853-20-9

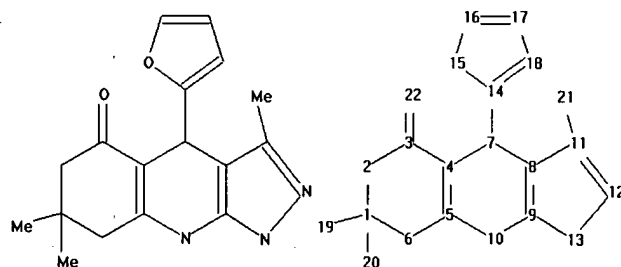
TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for
details.

Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>
Uploading H:\STN queries\10612885b.str



chain nodes :
19 20 21 22
ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18
chain bonds :
1-19 1-20 3-22 7-14 11-21
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10 9-13
11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact/norm bonds :
1-2 1-6 2-3 3-4 3-22 4-5 4-7 5-6 5-10 7-8 8-9 8-11 9-10
9-13 11-12 12-13 14-15 14-18 15-16 16-17 17-18
exact bonds :
1-19 1-20 7-14 11-21

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom
9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom
17:Atom 18:Atom 19:CLASS 20:CLASS 21:CLASS 22:CLASS

L1 STRUCTURE UPLOADED

=> d l1
L1 HAS NO ANSWERS
L1 STR

Page 1-A



Page 2-A

NODE ATTRIBUTES:
HCOUNT IS M3 AT 19
HCOUNT IS M3 AT 20
HCOUNT IS M3 AT 21
NSPEC IS R AT 1
NSPEC IS R AT 2
NSPEC IS R AT 3
NSPEC IS R AT 4
NSPEC IS R AT 5
NSPEC IS R AT 6
NSPEC IS R AT 7
NSPEC IS R AT 8
NSPEC IS R AT 9
NSPEC IS R AT 10
NSPEC IS R AT 11
NSPEC IS R AT 12
NSPEC IS R AT 13
NSPEC IS R AT 14
NSPEC IS R AT 15
NSPEC IS R AT 16
NSPEC IS R AT 17
NSPEC IS R AT 18
NSPEC IS C AT 19
NSPEC IS C AT 20

NSPEC IS C AT 21
NSPEC IS C AT 22
DEFAULT MLEVEL IS ATOM
MLEVEL IS CLASS AT 19 20 21 22
DEFAULT ELEVEL IS LIMITED

GRAPH ATTRIBUTES:
RING(S) ARE ISOLATED OR EMBEDDED
NUMBER OF NODES IS 22

STEREO ATTRIBUTES: NONE

=> s l1 sam
SAMPLE SEARCH INITIATED 10:25:56 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 0 TO 0
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s l1 fam
SAMPLE SEARCH INITIATED 10:26:03 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0
ANSWERS
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 0 TO 0
PROJECTED ANSWERS: 0 TO 0

L3 0 SEA FAM SAM L1

=> s l1 fu1
FULL SEARCH INITIATED 10:26:08 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 21 TO ITERATE

100.0% PROCESSED 21 ITERATIONS 8
ANSWERS
SEARCH TIME: 00.00.01

L4 8 SEA SSS FUL L1

=> s l1 fu1 fam
FULL SEARCH INITIATED 10:26:12 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 5 TO ITERATE

| | | | |
|------------------------------|--------------------|-------------|---------|
| Bioconc. Factor (BCF) | 98.8 | pH 8 | (1) ACD |
| Bioconc. Factor (BCF) | 98.8 | pH 10 | (1) ACD |
| Boiling Point (BP) | 478.2+/-45.0 deg C | 760 Torr | (1) ACD |
| Enthalpy of Vap. (HVP) | 74.23+/-3.0 kJ/mol | | (1) ACD |
| Flash Point (FP) | 243.0+/-51.7 deg C | | (1) ACD |
| Freely Rotatable Bonds (FRB) | 1 | | (1) ACD |
| H acceptors (HAC) | 5 | | (1) ACD |
| H donors (HD) | 2 | | (1) ACD |
| KOC (KOC) | 1 | pH 1 | (1) ACD |
| KOC (KOC) | 79.0 | pH 4 | (1) ACD |
| KOC (KOC) | 922 | pH 7 | (1) ACD |
| KOC (KOC) | 931 | pH 8 | (1) ACD |
| KOC (KOC) | 932 | pH 10 | (1) ACD |
| logD (LOGD) | -0.49 | pH 1 | (1) ACD |
| logD (LOGD) | 1.86 | pH 4 | (1) ACD |
| logD (LOGD) | 2.92 | pH 7 | (1) ACD |
| logD (LOGD) | 2.93 | pH 8 | (1) ACD |
| logD (LOGD) | 2.93 | pH 10 | (1) ACD |
| logP (LOGP) | 2.928+/-0.412 | | (1) ACD |
| Molar Solubility (SLB.MOL) | >=0.1 - <1 mol/L | pH 1 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 4 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 7 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 8 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 10 | (1) ACD |
| Molecular Weight (MW) | 297.35 | | (1) ACD |
| pKa (PKA) | 13.94+/-0.60 | Most Acidic | (1) ACD |
| pKa (PKA) | 4.82+/-0.60 | Most Basic | (1) ACD |
| Vapor Pressure (VP) | 2.62E-09 Torr | 25 deg C | (1) ACD |

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
Solaris v4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

REFERENCE 1

AN 140:87744 CA Full-text
TI Affinity small molecules for the EPO receptor
IN Olsson, Lennart; Naranda, Tatjana
PA Receptron, Inc., USA
SO PCT Int. Appl., 85 pp.
CODEN: PIXXD2
DT Patent
LA English
IC ICM C07K
CC 1-12 (Pharmacology)
Section cross-reference(s): 2

| FAN.CNT | PATENT NO. | KIND | DATE | APPLICATION NO. | DATE |
|---------|---------------|------|----------|-----------------|----------|
| PI | WO 2004005323 | A2 | 20040115 | WO 2003-US21394 | 20030703 |
| | WO 2004005323 | A3 | 20040701 | | |

100.0% PROCESSED 5 ITERATIONS 1
ANSWERS
SEARCH TIME: 00.00.01

L5 1 SEA FAM FUL L1

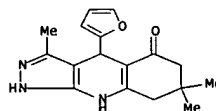
=> 15 not 14
L6 0 L5 NOT L4

=> d 15 1 all

L5 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2004 ACS on STM
RN 645337-25-3 REGISTRY
ED Entered STN: 03 Feb 2004
CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl- (9CI) (CA INDEX NAME)
FS 3D CONCORD
MF C17 H19 N3 O2
SR CA
LC STN Files: CA, CAPLUS, USPATFULL
DT.CA Caplus document type: Patent
RL.P Roles from patents: BIOL (Biological study); USES (Uses)

Ring System Data

| Elemental Analysis EA | Elemental Sequence ES | Size of the Rings SZ | Ring System Formula RF | Ring Identifier RID | RID Occurrence Count |
|-----------------------|-----------------------|----------------------|------------------------|---------------------|----------------------|
| C40 | 10C4 | 5 | C40 | 16.138.5 | 11 |
| C3N2-C5N-C6 | N2C3-NC5-C6 | 5-6-6 | C10N3 | 1894.85.24 | 1 |



Calculated Properties (CALC)

| PROPERTY (CODE) | VALUE | CONDITION | NOTE |
|-----------------------|-------|-----------|---------|
| Bioconc. Factor (BCF) | 1 | pH 1 | (1) ACD |
| Bioconc. Factor (BCF) | 8.39 | pH 4 | (1) ACD |
| Bioconc. Factor (BCF) | 97.8 | pH 7 | (1) ACD |

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZM, ZW
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG
US 2004171541 A1 20040902 US 2003-613754 20030702
US 2004116346 A1 20040617 US 2003-612885 20030703
PRAI US 2002-393360P 20020703
US 2002-393361P 20020703
US 2002-394110P 20020703
AB Comps. are provided that complex with the modulating domain of erythropoietin receptor (EPO-R) for use with EPO-R to determine the presence of EPO-R, the ability of other mols. to bind to the modulating domain in competitive assays and to induce a signal by EPO-R into a cell when bound by the subject compds. in a physiol. environment. The compds. are characterized by having a six-membered heterocyclic ring comprising at least one nitrogen atom and include substituted triazopyrimidine, pyridazinone, pyridine and piperidine.
ST EPO receptor modulator small mol
IT Proteins
RL: BSU (Biological study, unclassified); BIOL (Biological study) (Bcl-xL, expression; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT Peptides, biological studies
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(EPO receptor modulating sequence; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)
IT Cell membrane
(EPO receptors of; affinity small mols. for erythropoietin

(EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Anemia (disease)
Cell proliferation
Combinatorial library
Drug delivery systems
Drug screening
Erythrocyte
Erythropoiesis
Hematocrit
Hematopoietic precursor cell
Human
Reticulocyte
(affinity small mols. for erythropoietin (EPO) receptor and

EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Erythropoietin receptors
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and

EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Nerve
(neuron; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT Cytoprotective agents
(neuroprotective; affinity small mols. for erythropoietin (EPO) receptor and EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 2503-56-2 40775-78-8 51646-16-3 51646-17-4 51646-19-6 51646-43-6
56347-20-7 63901-48-4 90559-98-1 90815-61-5 113967-71-8
113967-74-1 194342-06-8 212074-47-0 244167-89-3 245082-87-5
245413-82-5 259683-29-9 261704-08-9 261704-09-0 262291-81-6

263267-38-5 287728-46-5 303145-64-4 303145-73-5 338793-16-1
645337-19-5 645337-20-8 645337-21-9 645337-22-0 645337-23-1
645337-24-2 645337-25-3
RL: BSU (Biological study, unclassified); BUU (Biological use, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and

EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 11096-26-7, Erythropoietin
RL: BSU (Biological study, unclassified); PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(affinity small mols. for erythropoietin (EPO) receptor and

EPO receptor modulating sequence in relation to modulating the response to the stimulus of hematopoietic or neuronal cells and treatment of anemia)

IT 239133-03-0 645415-22-1
RL: PRP (Properties)
(unclassified sequence; affinity small mols. for the EPO receptor)

=> d 14 1-8 ful
'FUL' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN
SAM - Index Name, MF, and structure - no RN
FIDE - All substance data, except sequence data
IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data
SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used
SQD - Protein sequence data, includes RN
SQD3 - Same as SQD, but 3-letter amino acid codes are used
SQN - Protein sequence name information, includes RN

CALC - Table of calculated properties
EPROP - Table of experimental properties
PROPP - EPROP and CALC

Any CA File format may be combined with any substance format to

obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract
APPS -- Application and Priority Information
BIB -- CA Accession Number, plus Bibliographic Data
CAN -- CA Accession Number
CBIB -- CA Accession Number, plus Bibliographic Data (compressed)
IND -- Index Data
IPC -- International Patent Classification
PATS -- PI, SO
STD -- BIB, IPC, and NCL

IABS -- ABS, indented, with text labels
IBIB -- BIB, indented, with text labels
ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)
OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.
The MAX format is the same as ALL.
The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

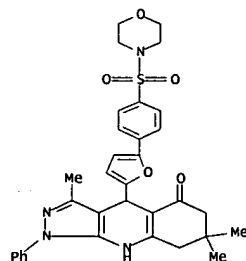
For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):all

L4 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
RN 748146-78-3 REGISTRY
ED Entered STN: 20 Sep 2004
CN INDEX NAME NOT YET ASSIGNED
FS 3D CONCORD
MF C33 H34 N4 O5 S
SR Chemical Library

Ring System Data

| Elemental Analysis EA | Elemental Sequence ES | Size of the Rings SZ | Ring System Formula RF | Ring Identifier RID | RID Occurrence Count |
|--------------------------|--------------------------|-------------------------|------------------------------|------------------------|----------------------|
| C4O | OC4 | 5 | C4O | 16.138.5 | 1 |
| C6 | C6 | 6 | C6 | 46.150.18 | 2 |
| C4NO | NC2OC2 | 6 | C4NO | 46.402.1 | 1 |
| C3N2-C5N-C6 | N2C3-NC5-C6 | 5-6-6 | C10N3 | 1894.85.24 | 1 |



Calculated Properties (CALC)

| PROPERTY (CODE) | VALUE | CONDITION | NOTE |
|------------------------------|---------------------|-----------|---------|
| Bioconc. Factor (BCF) | 12.13 | ph 1 | (1) ACD |
| Bioconc. Factor (BCF) | 649 | ph 4 | (1) ACD |
| Bioconc. Factor (BCF) | 1353 | ph 7 | (1) ACD |
| Bioconc. Factor (BCF) | 1354 | ph 8 | (1) ACD |
| Bioconc. Factor (BCF) | 1355 | ph 10 | (1) ACD |
| Boiling Point (BP) | 749.7+/-70.0 deg C | 760 Torr | (1) ACD |
| Enthalpy of Vap. (HVAP) | 109.28+/-3.0 kJ/mol | | (1) ACD |
| Flash Point (FP) | 407.2+/-64.2 deg C | | (1) ACD |
| Freely Rotatable Bonds (FRB) | 4 | | (1) ACD |
| H acceptors (HAC) | 9 | | (1) ACD |
| H donors (HD) | 1 | | (1) ACD |
| Koc (KOC) | 9.55 | ph 1 | (1) ACD |
| Koc (KOC) | 2910 | ph 4 | (1) ACD |
| Koc (KOC) | 6062 | ph 7 | (1) ACD |
| Koc (KOC) | 6067 | ph 8 | (1) ACD |
| Koc (KOC) | 6068 | ph 10 | (1) ACD |
| logD (LOGD) | 1.62 | ph 1 | (1) ACD |
| logD (LOGD) | 4.10 | ph 4 | (1) ACD |
| logD (LOGD) | 4.42 | ph 7 | (1) ACD |
| logD (LOGD) | 4.42 | ph 8 | (1) ACD |
| logD (LOGD) | 4.42 | ph 10 | (1) ACD |
| logD (LOGD) | 4.42+/-0.655 | | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | ph 1 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | ph 4 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | ph 7 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | ph 8 | (1) ACD |

Molar Solubility (SLB.MOL) <0.01 mol/L pH 10 (1) ACD
 Molecular Weight (MW) 598.71 (1) ACD
 pKa (PKA) 3.95+/-0.60 Most Basic (1) ACD
 Vapor Pressure (VP) 2.33E-22 Torr 25 deg C (1) ACD

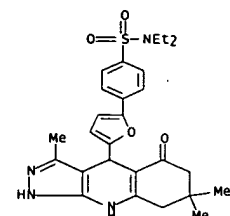
(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
 Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 748146-41-0 REGISTRY
 ED Entered STN: 20 Sep 2004
 CN INDEX NAME NOT YET ASSIGNED
 FS 3D CONCORD
 MF C27 H32 N4 O4 S
 SR Chemical Library

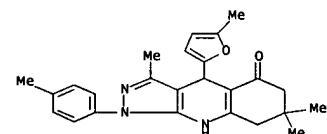
Ring System Data

| Elemental Analysis EA | Elemental Sequence ES | Size of the Rings SZ | Ring System Formula RF | Ring Identifier RID | RID Occurrence Count |
|-----------------------|-----------------------|----------------------|------------------------|---------------------|----------------------|
| C40 | OC4 | 5 | C40 | 16.138.5 | 1 |
| C6 | C6 | 6 | C6 | 46.150.18 | 1 |
| C3N2-C5N-C6 | N2C3-NC5-C6 | 5-6-6 | C10N3 | 1894.85.24 | 1 |



Calculated Properties (CALC)

C40 OC4 5 C40 16.138.5 1
 C6 C6 6 C6 46.150.18 1
 C3N2-C5N-C6 N2C3-NC5-C6 5-6-6 C10N3 1894.85.24 1



Calculated Properties (CALC)

| PROPERTY (CODE) | VALUE | CONDITION | NOTE |
|------------------------------|--------------------|------------|---------|
| Bioconc. Factor (BCF) | 2.22 | pH 1 | (1) ACD |
| Bioconc. Factor (BCF) | 847 | pH 4 | (1) ACD |
| Bioconc. Factor (BCF) | 2188 | pH 7 | (1) ACD |
| Bioconc. Factor (BCF) | 2191 | pH 8 | (1) ACD |
| Bioconc. Factor (BCF) | 2192 | pH 10 | (1) ACD |
| Boiling Point (BP) | 536.3+/-50.0 deg C | 760 Torr | (1) ACD |
| Enthalpy of Vap. (HVP) | 81.29+/-3.0 kJ/mol | | (1) ACD |
| Flash Point (FP) | 278.1+/-54.2 deg C | | (1) ACD |
| Freely Rotatable Bonds (FRB) | 2 | | (1) ACD |
| H acceptors (HAC) | 5 | | (1) ACD |
| H donors (HD) | 1 | | (1) ACD |
| Koc (KOC) | 8.69 | pH 1 | (1) ACD |
| Koc (KOC) | 3308 | pH 4 | (1) ACD |
| Koc (KOC) | 8549 | pH 7 | (1) ACD |
| Koc (KOC) | 8562 | pH 8 | (1) ACD |
| Koc (KOC) | 8563 | pH 10 | (1) ACD |
| logD (LOGD) | 1.70 | pH 1 | (1) ACD |
| logD (LOGD) | 4.29 | pH 4 | (1) ACD |
| logD (LOGD) | 4.70 | pH 7 | (1) ACD |
| logD (LOGD) | 4.70 | pH 8 | (1) ACD |
| logD (LOGD) | 4.70 | pH 10 | (1) ACD |
| logP (LOGP) | 4.699+/-0.511 | | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 1 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 4 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 7 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 8 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 10 | (1) ACD |
| Molecular Weight (MW) | 401.50 | | (1) ACD |
| pKa (PKA) | 4.16+/-0.60 | Most Basic | (1) ACD |
| Vapor Pressure (VP) | 1.42E-11 Torr | 25 deg C | (1) ACD |

| PROPERTY (CODE) | VALUE | CONDITION | NOTE |
|------------------------------|--------------------|-------------|---------|
| Bioconc. Factor (BCF) | 1.22 | pH 1 | (1) ACD |
| Bioconc. Factor (BCF) | 274 | pH 4 | (1) ACD |
| Bioconc. Factor (BCF) | 3190 | pH 7 | (1) ACD |
| Bioconc. Factor (BCF) | 3221 | pH 8 | (1) ACD |
| Bioconc. Factor (BCF) | 3225 | pH 10 | (1) ACD |
| Boiling Point (BP) | 680.7+/-65.0 deg C | 760 Torr | (1) ACD |
| Enthalpy of Vap. (HVP) | 99.88+/-3.0 kJ/mol | | (1) ACD |
| Flash Point (FP) | 365.5+/-61.7 deg C | | (1) ACD |
| Freely Rotatable Bonds (FRB) | 5 | | (1) ACD |
| H acceptors (HAC) | 8 | | (1) ACD |
| H donors (HD) | 2 | | (1) ACD |
| Koc (KOC) | 4.28 | pH 1 | (1) ACD |
| Koc (KOC) | 958 | pH 4 | (1) ACD |
| Koc (KOC) | 11170 | pH 7 | (1) ACD |
| Koc (KOC) | 11278 | pH 8 | (1) ACD |
| Koc (KOC) | 11289 | pH 10 | (1) ACD |
| logD (LOGD) | 1.50 | pH 1 | (1) ACD |
| logD (LOGD) | 3.85 | pH 4 | (1) ACD |
| logD (LOGD) | 4.91 | pH 7 | (1) ACD |
| logD (LOGD) | 4.92 | pH 8 | (1) ACD |
| logD (LOGD) | 4.92 | pH 10 | (1) ACD |
| logP (LOGP) | 4.920+/-0.532 | | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 1 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 4 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 7 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 8 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 10 | (1) ACD |
| Molecular Weight (MW) | 508.63 | | (1) ACD |
| pKa (PKA) | 13.94+/-0.60 | Most Acidic | (1) ACD |
| pKa (PKA) | 4.82+/-0.60 | Most Basic | (1) ACD |
| Vapor Pressure (VP) | 2.19E-18 Torr | 25 deg C | (1) ACD |

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
 Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 748145-15-5 REGISTRY
 ED Entered STN: 20 Sep 2004
 CN INDEX NAME NOT YET ASSIGNED
 FS 3D CONCORD
 MF C25 H27 N3 O2
 SR Chemical Library

Ring System Data

| Elemental Analysis EA | Elemental Sequence ES | Size of the Rings SZ | Ring System Formula RF | Ring Identifier RID | RID Occurrence Count |
|-----------------------|-----------------------|----------------------|------------------------|---------------------|----------------------|
|-----------------------|-----------------------|----------------------|------------------------|---------------------|----------------------|

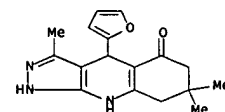
(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
 Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 645337-25-3 REGISTRY
 ED Entered STN: 03 Feb 2004
 CN 5H-pyrazolo[3,4-b]quinolin-5-one, 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-
 3,7,7-trimethyl- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C17 H19 N3 O2
 SR CA
 LC STN Files: CA, CAPLUS, USPATFULL
 DT.CA Caplus document type: Patent
 RL.P Roles from patents: BIOL (Biological study); USES (Uses)

Ring System Data

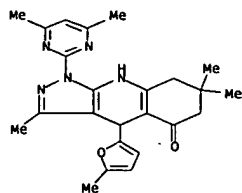
| Elemental Analysis EA | Elemental Sequence ES | Size of the Rings SZ | Ring System Formula RF | Ring Identifier RID | RID Occurrence Count |
|-----------------------|-----------------------|----------------------|------------------------|---------------------|----------------------|
| C40 | OC4 | 5 | C40 | 16.138.5 | 1 |
| C3N2-C5N-C6 | N2C3-NC5-C6 | 5-6-6 | C10N3 | 1894.85.24 | 1 |



Calculated Properties (CALC)

| PROPERTY (CODE) | VALUE | CONDITION | NOTE |
|-----------------------|--------------------|-----------|---------|
| Bioconc. Factor (BCF) | 1 | pH 1 | (1) ACD |
| Bioconc. Factor (BCF) | 8.39 | pH 4 | (1) ACD |
| Bioconc. Factor (BCF) | 97.8 | pH 7 | (1) ACD |
| Bioconc. Factor (BCF) | 98.8 | pH 8 | (1) ACD |
| Bioconc. Factor (BCF) | 98.8 | pH 10 | (1) ACD |
| Boiling Point (BP) | 478.2+/-45.0 deg C | 760 Torr | (1) ACD |

| Elemental Analysis EA | Elemental Sequence ES | Size of the Rings SZ | Ring System Formula RF | Ring Identifier RID | RID Occurrence Count |
|-----------------------|-----------------------|----------------------|------------------------|---------------------|----------------------|
| C4O | OC4 | 5 | C4O | 16.138.5 | 1 |
| C4N2 | NCNC3 | 6 | C4N2 | 46.195.39 | 1 |
| C3N2-C5N-C6 | C2C3-NC5-C6 | 5-6-6 | C10N3 | 1894.85.24 | 1 |



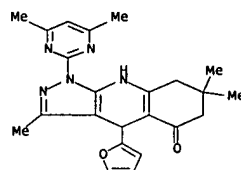
Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 521284-01-5 REGISTRY
 ED Entered STN: 28 May 2003
 CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 1-(4,6-dimethyl-2-pyrimidinyl)-4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C23 H25 N5 O2
 SR Chemical Library
 LC STN Files: CHEMCATS

Ring System Data

| Elemental Analysis EA | Elemental Sequence ES | Size of the Rings SZ | Ring System Formula RF | Ring Identifier RID | RID Occurrence Count |
|-----------------------|-----------------------|----------------------|------------------------|---------------------|----------------------|
| C40 | OC4 | 5 | C40 | 16.138.5 | 1 |
| C4N2 | NCNC3 | 6 | C4N2 | 46.195.39 | 1 |
| C3N2-C5N-C6 | N2C3-NC5-C6 | 5-6-6 | C10N3 | 1894.85.24 | 1 |



Calculated Properties (CALC)

| PROPERTY (CODE) | VALUE | CONDITION | NOTE |
|-----------------------|-------|-----------|---------|
| Bioconc. Factor (BCF) | 108 | pH 1 | (1) ACD |
| Bioconc. Factor (BCF) | 283 | pH 4 | (1) ACD |
| Bioconc. Factor (BCF) | 283 | pH 7 | (1) ACD |

Calculated Properties (CALC)

| PROPERTY (CODE) | VALUE | CONDITION | NOTE |
|------------------------------|--------------------|------------|---------|
| Bioconc. Factor (BCF) | 242 | pH 1 | (1) ACD |
| Bioconc. Factor (BCF) | 633 | pH 4 | (1) ACD |
| Bioconc. Factor (BCF) | 634 | pH 7 | (1) ACD |
| Bioconc. Factor (BCF) | 634 | pH 8 | (1) ACD |
| Bioconc. Factor (BCF) | 634 | pH 10 | (1) ACD |
| Boiling Point (BP) | 607.6+/-65.0 deg C | 760.0 Torr | (1) ACD |
| Enthalpy of Vap. (HVP) | 90.29+/-3.0 kJ/mol | | (1) ACD |
| Flash Point (FP) | 321.2+/-61.7 deg C | | (1) ACD |
| Freely Rotatable Bonds (FRB) | 2 | | (1) ACD |
| H acceptors (HAC) | 7 | | (1) ACD |
| H donors (HD) | 1 | | (1) ACD |
| Koc (KOC) | 1346 | pH 1 | (1) ACD |
| Koc (KOC) | 3518 | pH 4 | (1) ACD |
| Koc (KOC) | 3524 | pH 7 | (1) ACD |
| Koc (KOC) | 3524 | pH 8 | (1) ACD |
| Koc (KOC) | 3524 | pH 10 | (1) ACD |
| logD (LOGD) | 3.57 | pH 1 | (1) ACD |
| logD (LOGD) | 3.99 | pH 4 | (1) ACD |
| logD (LOGD) | 3.99 | pH 7 | (1) ACD |
| logD (LOGD) | 3.99 | pH 8 | (1) ACD |
| logD (LOGD) | 3.99 | pH 10 | (1) ACD |
| logP (LOGP) | 3.990+/-0.891 | | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 1 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 4 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 7 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 8 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 10 | (1) ACD |
| Molecular Weight (MW) | 417.50 | | (1) ACD |
| pKa (PKA) | 1.15+/-0.60 | Most Basic | (1) ACD |
| Vapor Pressure (VP) | 1.04E-14 Torr | 25.0 deg C | (1) ACD |

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software

| | | | |
|------------------------------|--------------------|------------|---------|
| Bioconc. Factor (BCF) | 283 | pH 8 | (1) ACD |
| Bioconc. Factor (BCF) | 283 | pH 10 | (1) ACD |
| Boiling Point (BP) | 601.5+/-65.0 deg C | 760.0 Torr | (1) ACD |
| Enthalpy of Vap. (HVP) | 89.51+/-3.0 kJ/mol | | (1) ACD |
| Flash Point (FP) | 317.6+/-61.7 deg C | | (1) ACD |
| Freely Rotatable Bonds (FRB) | 2 | | (1) ACD |
| H acceptors (HAC) | 7 | | (1) ACD |
| H donors (HD) | 1 | | (1) ACD |
| Koc (KOC) | 756 | pH 1 | (1) ACD |
| Koc (KOC) | 1977 | pH 4 | (1) ACD |
| Koc (KOC) | 1980 | pH 7 | (1) ACD |
| Koc (KOC) | 1980 | pH 8 | (1) ACD |
| Koc (KOC) | 1980 | pH 10 | (1) ACD |
| logD (LOGD) | 3.11 | pH 1 | (1) ACD |
| logD (LOGD) | 3.53 | pH 4 | (1) ACD |
| logD (LOGD) | 3.53 | pH 7 | (1) ACD |
| logD (LOGD) | 3.53 | pH 8 | (1) ACD |
| logD (LOGD) | 3.53 | pH 10 | (1) ACD |
| logP (LOGP) | 3.530+/-0.890 | | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 1 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 4 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 7 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 8 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 10 | (1) ACD |
| Molecular Weight (MW) | 403.48 | | (1) ACD |
| pKa (PKA) | 1.15+/-0.60 | Most Basic | (1) ACD |
| Vapor Pressure (VP) | 2.00E-14 Torr | 25.0 deg C | (1) ACD |

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software

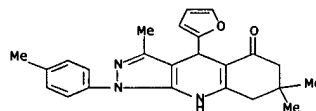
Solaris V4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

L4 ANSWER 7 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 380450-98-6 REGISTRY
 ED Entered STN: 04 Jan 2002
 CN 5H-Pyrazolo[3,4-b]quinolin-5-one, 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-trimethyl-1-(4-methylphenyl)- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C24 H25 N3 O2
 SR Chemical Library
 LC STN Files: CHEMCATS

Ring System Data

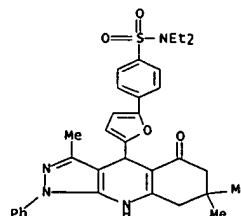
| Elemental Analysis EA | Elemental Sequence ES | Size of the Rings SZ | Ring System Formula RF | Ring Identifier RID | RID Occurrence Count |
|-----------------------|-----------------------|----------------------|------------------------|---------------------|----------------------|
| C40 | OC4 | 5 | C40 | 16.138.5 | 1 |
| C6 | C6 | 6 | C6 | 46.150.18 | 1 |
| C3N2-C5N-C6 | N2C3-NC5-C6 | 5-6-6 | C10N3 | 1894.85.24 | 1 |



L4 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
 RN 378189-53-8 REGISTRY
 ED Entered STN: 26 Dec 2001
 CN Benzenesulfonamide, N,N-diethyl-4-[5-(4,5,6,7,8,9-hexahydro-3,7,7-trimethyl-5-oxo-1-phenyl-1H-pyrazolo[3,4-b]quinolin-4-yl)-2-furanyl]- (9CI) (CA INDEX NAME)
 FS 3D CONCORD
 MF C33 H36 N4 O4 S
 SR Chemical Library
 LC STN Files: CHEMCATS

Ring System Data

| Elemental Analysis EA | Elemental Sequence ES | Size of the Rings SZ | Ring System Formula RF | Ring Identifier RID | RID Occurrence Count |
|-----------------------|-----------------------|----------------------|------------------------|---------------------|----------------------|
| C40 | OC4 | 5 | C40 | 16.138.5 | 1 |
| C6 | C6 | 6 | C6 | 46.150.18 | 2 |
| C3N2-C5N-C6 | N2C3-NC5-C6 | 5-6-6 | C10N3 | 1894.85.24 | 1 |



Calculated Properties (CALC)

| PROPERTY (CODE) | VALUE | CONDITION | NOTE |
|------------------------------|---------------------|------------|---------|
| Bioconc. Factor (BCF) | 22.5 | pH 1 | (1) ACD |
| Bioconc. Factor (BCF) | 6852 | pH 4 | (1) ACD |
| Bioconc. Factor (BCF) | 14276 | pH 7 | (1) ACD |
| Bioconc. Factor (BCF) | 14290 | pH 8 | (1) ACD |
| Bioconc. Factor (BCF) | 14291 | pH 10 | (1) ACD |
| Boiling Point (BP) | 712.3+/-70.0 deg C | 760.0 Torr | (1) ACD |
| Enthalpy of Vap. (HVP) | 104.14+/-3.0 kJ/mol | | (1) ACD |
| Flash Point (FP) | 384.6+/-64.2 deg C | | (1) ACD |
| Freely Rotatable Bonds (FRB) | 6 | | (1) ACD |
| H acceptors (HAC) | 8 | | (1) ACD |
| H donors (HD) | 1 | | (1) ACD |
| Koc (KOC) | 51.6 | pH 1 | (1) ACD |
| Koc (KOC) | 15714 | pH 4 | (1) ACD |
| Koc (KOC) | 32737 | pH 7 | (1) ACD |
| Koc (KOC) | 32769 | pH 8 | (1) ACD |
| Koc (KOC) | 32772 | pH 10 | (1) ACD |
| logD (LOGD) | 2.97 | pH 1 | (1) ACD |
| logD (LOGD) | 5.45 | pH 4 | (1) ACD |
| logD (LOGD) | 5.77 | pH 7 | (1) ACD |
| logD (LOGD) | 5.77 | pH 8 | (1) ACD |
| logD (LOGD) | 5.77 | pH 10 | (1) ACD |
| logP (LOGP) | 5.771+/-0.615 | | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 1 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 4 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 7 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 8 | (1) ACD |
| Molar Solubility (SLB.MOL) | <0.01 mol/L | pH 10 | (1) ACD |
| Molecular Weight (Mw) | 584.73 | | (1) ACD |
| pKa (PKA) | 3.95+/-0.60 | Most Basic | (1) ACD |
| Vapor Pressure (VP) | 3.87E-20 Torr | 25.0 deg C | (1) ACD |

(1) Calculated using Advanced Chemistry Development (ACD/Labs) Software
Solaris v4.76 ((C) 1994-2004 ACD/Labs)

See HELP PROPERTIES for information about property data sources in REGISTRY.

=> d 14 1-8 ed

L4 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 20 Sep 2004

L4 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 20 Sep 2004

FILE 'HCAPLUS' ENTERED AT 10:35:54 ON 28 DEC 2004
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=> d his

(FILE 'HOME' ENTERED AT 10:25:10 ON 28 DEC 2004)

FILE 'REGISTRY' ENTERED AT 10:25:22 ON 28 DEC 2004
STRUCTURE UPLOADED
L1 0 S L1 SAM
L2 0 S L1 FAM
L3 8 S L1 FUL
L4 1 S L1 FUL FAM
L5 0 L5 NOT L4

FILE 'HCAPLUS, USPATFULL, CHEMCATS' ENTERED AT 10:35:54 ON 28 DEC 2004

=> s 14
L7 9 L4

=> d 19 ed
L9 NOT FOUND

The L-number entered has not been defined in this session, or it has been deleted. To see the L-numbers currently defined in this session, enter DISPLAY HISTORY at an arrow prompt (=>).

=> d 17 1-9 ed
NO VALID FORMATS ENTERED FOR FILE 'USPATFULL'
NO VALID FORMATS ENTERED FOR FILE 'CHEMCATS'
In a multifile environment, each file must have at least one valid format requested. Refer to file specific help messages or the STNGUIDE file for information on formats available in individual files.
REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):end

=> d 17 1-9 ibib abs
NO VALID FORMATS ENTERED FOR FILE 'CHEMCATS'
In a multifile environment, each file must have at least one valid format requested. Refer to file specific help messages or the STNGUIDE file for information on formats available in individual files.
REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):
REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):filedefault

L7 ANSWER 1 OF 9 HCAPLUS COPYRIGHT 2004 ACS on STN

L4 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 20 Sep 2004

L4 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 03 Feb 2004

L4 ANSWER 5 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 28 May 2003

L4 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 28 May 2003

L4 ANSWER 7 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 04 Jan 2002

L4 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
ED Entered STN: 26 Dec 2001

=> d 14 1-8 1c

L4 ANSWER 1 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

L4 ANSWER 2 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

L4 ANSWER 3 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN

L4 ANSWER 4 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
LC STN Files: CA, CAPLUS, USPATFULL

L4 ANSWER 5 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
LC STN Files: CHEMCATS

L4 ANSWER 6 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
LC STN Files: CHEMCATS

L4 ANSWER 7 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
LC STN Files: CHEMCATS

L4 ANSWER 8 OF 8 REGISTRY COPYRIGHT 2004 ACS on STN
LC STN Files: CHEMCATS

=> fil hcaplus uspatfull chemcats

COST IN U.S. DOLLARS SINCE FILE

TOTAL ENTRY

SESSION FULL ESTIMATED COST 294.47
294.68

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE

TOTAL ENTRY

SESSION CA SUBSCRIBER PRICE -1.32
1.32

AN 2004:41501 HCAPLUS Full-text

DN 140:87744

TI Affinity small molecules for the EPO receptor

IN Olsson, Lennart; Naranda, Tatjana

PA Reception, Inc., USA

SO PCT Int. Appl., 85 pp.

CODEN: PIXXD2

DT Patent

LA English

FAN.CNT 1

PATENT NO. KIND DATE APPLICATION NO.

DATE -----

PI WO 2004005323 A2 20040115 WO 2003-US21394

20030703 WO 2004005323 A3 20040701

W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA,

CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD,

GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC,

LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ,

OM, PH, PL, PT, RO, RU, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT,

TZ, UA, UG, UZ, VN, YU, ZA, ZM, ZW

RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM,

EE, ES, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK,

SK, TR, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI,

TD, TG, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN,

US 2004171541 A1 20040902 US 2003-613754

20030702 US 2004116346 A1 20040617 US 2003-612885

20030703

PRAI US 2002-393360P P 20020703

US 2002-393361P P 20020703

US 2002-394110P P 20020703

OS MARPAT 140:87744

L7 ANSWER 2 OF 9 USPATFULL on STN

AN 2004:221770 USPATFULL Full-text

TI Affinity small molecules for the EPO receptor

IN Olsson, Lennart, Orinda, CA, UNITED STATES

Naranda, Tatjana, Mountain View, CA, UNITED STATES

PI US 2004171541 A1 20040902

AI US 2003-613754 A1 20030702 (10)

PRAI US 2002-393360P 20020703 (60)

US 2002-393361P 20020703 (60)

US 2002-394110P 20020703 (60)

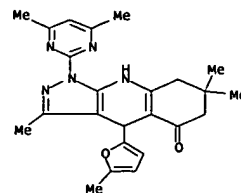
DT Utility

FS APPLICATION

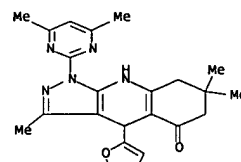
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 NCL NCLM: 514/012.000
 514/348.000 NCL: 514/247.000; 514/259.310; 514/292.000; 514/347.000;
 IC [7]
 ICM: A61K038-18
 ICS: A61K031-50; A61K031-519; A61K031-44; A61K031-4745
 CAS INDEXING IS AVAILABLE FOR THIS PATENT.

L7 ANSWER 3 OF 9 USPATFULL on STN
 AN 2004:152124 USPATFULL Full-text
 TI Affinity small molecules for the EPO receptor
 IN Olsson, Lennart, Orinda, CA, UNITED STATES
 Naranda, Tatjana, Mountain View, CA, UNITED STATES
 PI US 2004116346 A1 20040617
 AI US 2003-612885 A1 20030703 (10)
 PRAI US 2002-393361P 20020703 (60)
 US 2002-393360P 20020703 (60)
 US 2002-394110P 20020703 (60)
 DT Utility
 FS APPLICATION
 LN.CNT 2000
 INCL INCLM: 514/012.000
 514/348.000 INCL: 514/247.000; 514/259.310; 514/292.000; 514/347.000;
 NCL NCLM: 514/012.000
 514/348.000 NCL: 514/247.000; 514/259.310; 514/292.000; 514/347.000;
 IC [7]
 ICM: A61K038-17
 ICS: A61K031-519; A61K031-501; A61K031-4745; A61K031-44
 CAS INDEXING IS AVAILABLE FOR THIS PATENT.

L7 ANSWER 4 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:3262248 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0509-4266
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 1-(4,6-dimethyl-2-pyrimidinyl)-1,4,6,7,8,9-
 hexahydro-
 3,7,7-trimethyl-4-(5-methyl-2-furanyl)-
 CAS Registry No. (RN): 521318-71-8
 Supplementary Term (ST): CHEMICAL LIBRARY
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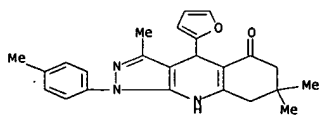


L7 ANSWER 5 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:3262236 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0509-4092
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 1-(4,6-dimethyl-2-pyrimidinyl)-4-(2-
 furanyl)-
 1,4,6,7,8,9-hexahydro-3,7,7-trimethyl-
 CAS Registry No. (RN): 521284-01-5
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :

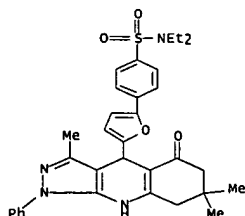


L7 ANSWER 6 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:2881171 CHEMCATS
 Catalog Name (CO): Enamine Screening Library
 Publication Date (PD): 30 Jun 2004
 Order Number (ON): T0508-1321
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-
 trimethyl-1-
 (4-methylphenyl)-
 CAS Registry No. (RN): 380450-98-6
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :

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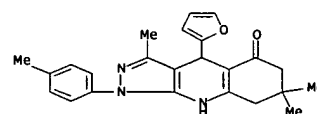


L7 ANSWER 7 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2003:2877484 CHEMCATS
 Catalog Name (CO): Enamine Screening Library
 Publication Date (PD): 30 Jun 2004
 Order Number (ON): T0505-5972
 Chemical Name (CN): Benzenesulfonamide, N,N-diethyl-4-[5-
 (4,5,6,7,8,9-
 hexahydro-3,7,7-trimethyl-5-oxo-1-phenyl]-
 1H-
 pyrazolo[3,4-b]quinolin-4-yl]-2-furanyl]-
 CAS Registry No. (RN): 378189-53-8
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :

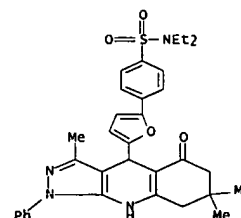


L7 ANSWER 8 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2002:1507742 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0508-1321
 Chemical Name (CN): 5H-Pyrazolo[3,4-b]quinolin-5-one,
 4-(2-furanyl)-1,4,6,7,8,9-hexahydro-3,7,7-
 trimethyl-1-
 (4-methylphenyl)-

CAS Registry No. (RN): 380450-98-6
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :



L7 ANSWER 9 OF 9 CHEMCATS COPYRIGHT 2004 ACS on STN
 Accession No. (AN): 2002:1491648 CHEMCATS
 Catalog Name (CO): Ambinter Screening Library
 Publication Date (PD): 1 Jan 2004
 Order Number (ON): T0505-5972
 Chemical Name (CN): Benzenesulfonamide, N,N-diethyl-4-[5-
 (4,5,6,7,8,9-
 hexahydro-3,7,7-trimethyl-5-oxo-1-phenyl]-
 1H-
 pyrazolo[3,4-b]quinolin-4-yl]-2-furanyl]-
 CAS Registry No. (RN): 378189-53-8
 Supplementary Term (ST): CHEMICAL LIBRARY
 Structure :



=> d 17 1-9 pd
 NO VALID FORMATS ENTERED FOR FILE 'HCAPLUS'
 NO VALID FORMATS ENTERED FOR FILE 'USPATFULL'
 In a multi-file environment, each file must have at least one valid
 format requested. Refer to file specific help messages or the
 STNGUIDE file for information on formats available in individual

files.
REENTER DISPLAY FORMAT FOR ALL FILES (FILEDEFAULT):end

=> index biosci
FILE 'DRUGMONOG' ACCESS NOT AUTHORIZED
COST IN U.S. DOLLARS
TOTAL
SESSION
FULL ESTIMATED COST 19.21
313.89
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)
TOTAL
ENTRY
SESSION
CA SUBSCRIBER PRICE 0.00
1.32
INDEX 'ADISCTI, ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, ANTE, AQUALINE, AQUASCI, BIOBUSINESS, BIOCOMMERCE, BIOENG, BIOSIS, BIOTECHABS, BIOTECHDS, BIOTECHNO, CABA, CANCERLIT, CAPLUS, CEABA-VTB, CEN, CIN, CONFSCI, CROPB, CROPU, DDFB, DDFU, DGENE, DISSABS, ...' ENTERED AT 10:38:10 ON 28 DEC 2004

75 FILES IN THE FILE LIST IN STINDEX

Enter SET DETAIL ON to see search term postings or to view search error messages that display as 0* with SET DETAIL OFF.

=> s (epo or (erythropoietin receptor))

FILE 'ADISCTI'
68 EPO
3 EPOS
71 EPO
(EPO OR EPOS)
641 ERYTHROPOIETIN
1 ERYTHROPOIETINS
642 ERYTHROPOIETIN
(ERYTHROPOIETIN OR ERYTHROPOIETINS)
77336 RECEPTOR
5437 RECEPTORS
79811 RECEPTOR
(RECEPTOR OR RECEPTORS)
3 ERYTHROPOIETIN RECEPTOR
(ERYTHROPOIETIN(W)RECEPTOR)
73 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'ADISINSIGHT'
24 EPO
2 EPOS
25 EPO
(EPO OR EPOS)
57 "ERYTHROPOIETIN"
15 "ERYTHROPOIETINS"

(RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
(ERYTHROPOIETIN(W)RECEPTOR)
22 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'AQUALINE'
1 EPO
2 EPOS
3 EPO
(EPO OR EPOS)
1 ERYTHROPOIETIN
241 RECEPTOR
286 RECEPTORS
395 RECEPTOR
(RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
(ERYTHROPOIETIN(W)RECEPTOR)
3 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'AQUASCI'
36 EPO
63 EPOS
99 EPO
(EPO OR EPOS)
10 "ERYTHROPOIETIN"
5274 "RECEPTOR"
4864 "RECEPTORS"
7710 "RECEPTOR"

("RECEPTOR" OR "RECEPTORS")
0 ERYTHROPOIETIN RECEPTOR
(("ERYTHROPOIETIN"(W)"RECEPTOR"))
99 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'BIOBUSINESS'
387 EPO
18 EPOS
404 EPO
(EPO OR EPOS)
588 "ERYTHROPOIETIN"
2 "ERYTHROPOIETINS"
590 "ERYTHROPOIETIN"
(("ERYTHROPOIETIN" OR "ERYTHROPOIETINS"))
6807 "RECEPTOR"
2441 "RECEPTORS"
7936 "RECEPTOR"
(("RECEPTOR" OR "RECEPTORS"))
7 ERYTHROPOIETIN RECEPTOR
(("ERYTHROPOIETIN"(W)"RECEPTOR"))
410 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'BIOCOMMERCE'
1192 EPO
5 EPOS
1197 EPO
(EPO OR EPOS)
837 ERYTHROPOIETIN
1 ERYTHROPOIETINS
837 ERYTHROPOIETIN
(ERYTHROPOIETIN OR ERYTHROPOIETINS)
3060 RECEPTOR
1032 RECEPTORS

57 "ERYTHROPOIETIN"
(("ERYTHROPOIETIN" OR "ERYTHROPOIETINS"))
4128 "RECEPTOR"
1761 "RECEPTORS"
4546 "RECEPTOR"
(("RECEPTOR" OR "RECEPTORS"))
6 ERYTHROPOIETIN RECEPTOR
(("ERYTHROPOIETIN"(W)"RECEPTOR"))
27 (EPO OR (ERYTHROPOIETIN RECEPTOR))
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16 EPO
162 ERYTHROPOIETIN
5 ERYTHROPOIETINS
163 ERYTHROPOIETIN
(ERYTHROPOIETIN OR ERYTHROPOIETINS)
1962 RECEPTOR
556 RECEPTORS
2258 RECEPTOR
(RECEPTOR OR RECEPTORS)
3 ERYTHROPOIETIN RECEPTOR
(ERYTHROPOIETIN(W)RECEPTOR)
19 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'AGRICOLA'
35 EPO
26 EPOS
60 EPO
(EPO OR EPOS)
136 ERYTHROPOIETIN
8600 RECEPTOR
9839 RECEPTORS
13992 RECEPTOR
(RECEPTOR OR RECEPTORS)
3 ERYTHROPOIETIN RECEPTOR
(ERYTHROPOIETIN(W)RECEPTOR)
63 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'ANABSTR'
26 EPO
27 EPOS
52 EPO
(EPO OR EPOS)
72 ERYTHROPOIETIN
1181 RECEPTOR
381 RECEPTORS
1391 RECEPTOR
(RECEPTOR OR RECEPTORS)
1 ERYTHROPOIETIN RECEPTOR
(ERYTHROPOIETIN(W)RECEPTOR)
53 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'ANTE'
10 EPO
12 EPOS
22 EPO
(EPO OR EPOS)
7 ERYTHROPOIETIN
91 RECEPTOR
73 RECEPTORS
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3917 RECEPTOR
(RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
(ERYTHROPOIETIN(W)RECEPTOR)
1197 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'BIOENG'
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183 EPO
(EPO OR EPOS)
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9 ERYTHROPOIETINS
336 ERYTHROPOIETIN
(ERYTHROPOIETIN OR ERYTHROPOIETINS)
11378 RECEPTOR
8432 RECEPTORS
14157 RECEPTOR
(RECEPTOR OR RECEPTORS)
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(ERYTHROPOIETIN(W)RECEPTOR)
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5822 EPO
(EPO OR EPOS)
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64 ERYTHROPOIETINS
17897 ERYTHROPOIETIN
(ERYTHROPOIETIN OR ERYTHROPOIETINS)
651542 RECEPTOR
323743 RECEPTORS
780807 RECEPTOR
(RECEPTOR OR RECEPTORS)
1356 ERYTHROPOIETIN RECEPTOR
(ERYTHROPOIETIN(W)RECEPTOR)
6546 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'BIOTECHABS'
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8 EPOS
356 EPO
(EPO OR EPOS)
981 ERYTHROPOIETIN
6 ERYTHROPOIETINS
983 ERYTHROPOIETIN
(ERYTHROPOIETIN OR ERYTHROPOIETINS)
16255 RECEPTOR
4183 RECEPTORS
17462 RECEPTOR
(RECEPTOR OR RECEPTORS)
85 ERYTHROPOIETIN RECEPTOR
(ERYTHROPOIETIN(W)RECEPTOR)
407 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'BIOTECHDS'
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8 EPOS
356 EPO

(EPO OR EPOS)
 981 ERYTHROPOIETIN
 6 ERYTHROPOIETINS
 983 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 16255 RECEPTOR
 4183 RECEPTORS
 17462 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 85 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 407 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'BIOTECHNO'
 2598 EPO
 33 EPOS
 2614 EPO
 (EPO OR EPOS)
 9222 ERYTHROPOIETIN
 29 ERYTHROPOIETINS
 9222 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 202158 RECEPTOR
 78475 RECEPTORS
 213334 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 774 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 2960 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CABA'
 247 EPO
 47 EPOS
 294 EPO
 (EPO OR EPOS)
 539 ERYTHROPOIETIN
 30692 RECEPTOR
 27815 RECEPTORS
 42614 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 21 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 308 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CANCERLIT'
 3530 EPO
 31 EPOS
 3544 EPO
 (EPO OR EPOS)
 11104 ERYTHROPOIETIN
 30 ERYTHROPOIETINS
 11104 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 122617 RECEPTOR
 117416 RECEPTORS
 160576 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 501 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 3775 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 539 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CONFSCI'
 87 EPO
 10 EPOS
 97 EPO
 (EPO OR EPOS)
 525 "ERYTHROPOIETIN"
 2 "ERYTHROPOIETINS"
 527 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 17313 "RECEPTOR"
 8593 "RECEPTORS"
 25653 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 71 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")
 165 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CROPB'
 0 EPO
 9 EPOS
 9 EPO
 (EPO OR EPOS)
 0 ERYTHROPOIETIN
 160 RECEPTOR
 35 RECEPTORS
 175 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 9 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CROPU'
 55 EPO
 12 EPOS
 67 EPO
 (EPO OR EPOS)
 1 ERYTHROPOIETIN
 1368 RECEPTOR
 570 RECEPTORS
 1566 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 67 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DDFB'
 2 EPO
 843 ERYTHROPOIETIN
 15280 RECEPTOR
 4633 RECEPTORS
 16551 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 3 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DDFU'
 1419 EPO
 3 EPOS
 1421 EPO

FILE 'CAPLUS'
 5521 EPO
 131 EPOS
 5625 EPO
 (EPO OR EPOS)
 11125 ERYTHROPOIETIN
 520 ERYTHROPOIETINS
 11155 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 582865 RECEPTOR
 534615 RECEPTORS
 693859 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1237 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 6216 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CEABA-VTB'
 106 EPO
 16 EPOS
 122 EPO
 (EPO OR EPOS)
 191 ERYTHROPOIETIN
 9 ERYTHROPOIETINS
 196 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 2128 RECEPTOR
 850 RECEPTORS
 2456 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 4 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 124 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CEN'
 33 EPO
 67 "ERYTHROPOIETIN"
 439 "RECEPTOR"
 381 "RECEPTORS"
 630 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 1 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")
 34 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'CIN'
 533 EPO
 6 EPOS
 536 EPO
 (EPO OR EPOS)
 544 "ERYTHROPOIETIN"
 4 "ERYTHROPOIETINS"
 547 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 4037 "RECEPTOR"
 1359 "RECEPTORS"
 4824 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 7 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")
 (EPO OR EPOS)
 3133 ERYTHROPOIETIN
 9 ERYTHROPOIETINS
 3133 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 99698 RECEPTOR
 44104 RECEPTORS
 111042 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 56 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 1441 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DGENE'
 15416 EPO
 68 EPOS
 15418 EPO
 (EPO OR EPOS)
 18662 ERYTHROPOIETIN
 21 ERYTHROPOIETINS
 18680 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 496912 RECEPTOR
 170378 RECEPTORS
 537954 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1551 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 15707 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DISSABS'
 152 EPO
 51 EPOS
 200 EPO
 (EPO OR EPOS)
 240 ERYTHROPOIETIN
 22949 RECEPTOR
 13830 RECEPTORS
 27515 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 35 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 216 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DRUGB'
 2 EPO
 843 ERYTHROPOIETIN
 15280 RECEPTOR
 4633 RECEPTORS
 16551 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 1 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 3 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DRUGMONOGZ'
 74 EPO
 2 EPOS
 76 EPO
 (EPO OR EPOS)
 1 ERYTHROPOIETIN

0 RECEPTOR
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 76 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'DRUGU'
 1633 EPO
 7 EPOS
 1637 EPO
 (EPO OR EPOS)
 3396 ERYTHROPOIETIN
 10 ERYTHROPOIETINS
 3396 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 110636 RECEPTOR
 54912 RECEPTORS
 125028 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 64 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 1657 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'EMBAL'
 52 EPO
 2 EPOS
 54 EPO
 (EPO OR EPOS)
 106 ERYTHROPOIETIN
 2 ERYTHROPOIETINS
 108 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 5008 RECEPTOR
 2397 RECEPTORS
 5899 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 6 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 57 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'EMBASE'
 4482 EPO
 111 EPOS
 4572 EPO
 (EPO OR EPOS)
 18291 "ERYTHROPOIETIN"
 51 "ERYTHROPOIETINS"
 18292 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 714332 "RECEPTOR"
 260226 "RECEPTORS"
 755991 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 1098 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)"RECEPTOR")
 5095 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'ESBIOBASE'
 1925 EPO
 44 EPOS
 1960 EPO
 (EPO OR EPOS)

FILE 'FSTA'
 21 EPO
 5 EPOS
 25 EPO
 (EPO OR EPOS)
 1 ERYTHROPOIETIN
 1012 RECEPTOR
 434 RECEPTORS
 1297 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 25 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'GENBANK'
 6183 EPO
 6652 "ERYTHROPOIETIN"
 201129 "RECEPTOR"
 289 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)"RECEPTOR")
 6457 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'HEALSAFE'
 7 EPO
 19 "ERYTHROPOIETIN"
 460 "RECEPTOR"
 311 "RECEPTORS"
 657 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 0 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)"RECEPTOR")
 7 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'IFIPAT'
 716 EPO
 11 EPOS
 725 EPO
 (EPO OR EPOS)
 1216 ERYTHROPOIETIN
 24 ERYTHROPOIETINS
 1235 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 32172 RECEPTOR
 11668 RECEPTORS
 36258 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 68 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 758 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'IMSDRUGNEWS'
 70 EPO
 161 "ERYTHROPOIETIN"
 3340 "RECEPTOR"
 836 "RECEPTORS"
 3873 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 7 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)"RECEPTOR")
 74 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'IMSPRODUCT'

3664 ERYTHROPOIETIN
 11 ERYTHROPOIETINS
 3666 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 239241 RECEPTOR
 129772 RECEPTORS
 280062 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 433 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 2150 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'FEDRIP'
 118 EPO
 2 EPOS
 120 EPO
 (EPO OR EPOS)
 243 ERYTHROPOIETIN
 16896 RECEPTOR
 9138 RECEPTORS
 19085 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 30 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 139 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'FOMAD'
 1 EPO
 12 EPOS
 13 EPO
 (EPO OR EPOS)
 1 ERYTHROPOIETIN
 2 RECEPTOR
 1 RECEPTORS
 2 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 13 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'FOREGE'
 0 EPO
 0 ERYTHROPOIETIN
 0 RECEPTOR
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 0 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'FROSTI'
 15 EPO
 15 EPOS
 29 EPO
 (EPO OR EPOS)
 9 ERYTHROPOIETIN
 1085 RECEPTOR
 784 RECEPTORS
 1582 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 29 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 31 EPO
 2 EPOS
 33 EPO
 (EPO OR EPOS)
 219 "ERYTHROPOIETIN"
 44 "RECEPTOR"
 8 "RECEPTORS"
 52 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 0 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)"RECEPTOR")
 33 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'IMSRESEARCH'
 32 EPO
 70 "ERYTHROPOIETIN"
 2213 "RECEPTOR"
 1048 "RECEPTORS"
 2816 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 6 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN(W)"RECEPTOR")
 36 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'JICST-EPLUS'
 621 EPO
 17 EPOS
 635 EPO
 (EPO OR EPOS)
 3631 ERYTHROPOIETIN
 4 ERYTHROPOIETINS
 3631 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 73555 RECEPTOR
 12647 RECEPTORS
 76150 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 192 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 800 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'KOSMET'
 5 EPO
 1 EPOS
 6 EPO
 (EPO OR EPOS)
 8 ERYTHROPOIETIN
 490 RECEPTOR
 555 RECEPTORS
 710 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 6 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'LIFESCI'
 863 EPO
 50 EPOS
 907 EPO
 (EPO OR EPOS)
 1458 "ERYTHROPOIETIN"

11 "ERYTHROPOIETINS"
 1458 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 167986 "RECEPTOR"
 153110 "RECEPTORS"
 216519 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 281 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")
 1047 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'MEDICINF'
 4 EPO
 7 EPOS
 11 EPO
 (EPO OR EPOS)
 22 ERYTHROPOIETIN
 1 ERYTHROPOIETINS
 23 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 324 RECEPTOR
 571 RECEPTORS
 827 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 11 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'MEDLINE'
 4856 EPO
 99 EPOS
 4929 EPO
 (EPO OR EPOS)
 16393 ERYTHROPOIETIN
 53 ERYTHROPOIETINS
 16397 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 479689 RECEPTOR
 514892 RECEPTORS
 673211 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 714 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 5281 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'NIOSHITIC'
 15 EPO
 89 ERYTHROPOIETIN
 1 ERYTHROPOIETINS
 89 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 1641 RECEPTOR
 1030 RECEPTORS
 2222 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 15 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'NTIS'
 46 EPO

694 RECEPTORS
 9243 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 0 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PHAR'
 72 EPO
 85 "ERYTHROPOIETIN"
 9931 "RECEPTOR"
 1111 "RECEPTORS"
 10104 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 47 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")
 73 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PHARMAML'
 99 EPO
 3 EPOS
 101 EPO
 (EPO OR EPOS)
 202 ERYTHROPOIETIN
 5 ERYTHROPOIETINS
 206 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 1768 RECEPTOR
 465 RECEPTORS
 2031 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 101 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PHIC'
 7 EPO
 6 "ERYTHROPOIETIN"
 18 "RECEPTOR"
 5 "RECEPTORS"
 23 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 0 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")
 7 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PHIN'
 841 EPO
 22 EPOS
 859 EPO
 (EPO OR EPOS)
 1142 "ERYTHROPOIETIN"
 18 "ERYTHROPOIETINS"
 1153 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 4441 "RECEPTOR"
 2186 "RECEPTORS"
 5782 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 3 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")

64 EPOS
 108 EPO
 (EPO OR EPOS)
 121 ERYTHROPOIETIN
 25 ERYTHROPOIETINS
 122 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 4927 RECEPTOR
 3307 RECEPTORS
 6501 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 108 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'NUTRACEUT'
 11 EPO
 0 ERYTHROPOIETIN
 10 RECEPTOR
 13 RECEPTORS
 22 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 11 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'OCEAN'
 24 EPO
 51 EPOS
 75 EPO
 (EPO OR EPOS)
 1 "ERYTHROPOIETIN"
 601 "RECEPTOR"
 581 "RECEPTORS"
 950 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 0 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")
 75 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PASCAL'
 2055 EPO
 127 EPOS
 2173 EPO
 (EPO OR EPOS)
 6424 ERYTHROPOIETIN
 22 ERYTHROPOIETINS
 6427 ERYTHROPOIETIN
 (ERYTHROPOIETIN OR ERYTHROPOIETINS)
 304437 RECEPTOR
 102112 RECEPTORS
 327238 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 334 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 2342 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PCTGEN'
 0 EPO
 0 ERYTHROPOIETIN
 8549 RECEPTOR

861 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PROMT'
 2713 EPO
 1902 EPOS
 4590 EPO
 (EPO OR EPOS)
 2025 "ERYTHROPOIETIN"
 34 "ERYTHROPOIETINS"
 2050 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 17544 "RECEPTOR"
 9487 "RECEPTORS"
 22972 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 27 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")
 4612 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PROUSDDR'
 22 EPO
 43 "ERYTHROPOIETIN"
 2 "ERYTHROPOIETINS"
 44 "ERYTHROPOIETIN"
 ("ERYTHROPOIETIN" OR "ERYTHROPOIETINS")
 36740 "RECEPTOR"
 19875 "RECEPTORS"
 44173 "RECEPTOR"
 ("RECEPTOR" OR "RECEPTORS")
 4 ERYTHROPOIETIN RECEPTOR
 ("ERYTHROPOIETIN"(W)"RECEPTOR")
 22 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'PS'
 0 EPO
 1 EPOS
 1 EPO
 (EPO OR EPOS)
 0 ERYTHROPOIETIN
 29 RECEPTOR
 1 RECEPTORS
 30 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 1 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'RDISCLOSURE'
 21 EPO
 2 EPOS
 23 EPO
 (EPO OR EPOS)
 1 ERYTHROPOIETIN
 93 RECEPTOR
 20 RECEPTORS
 106 RECEPTOR
 (RECEPTOR OR RECEPTORS)
 0 ERYTHROPOIETIN RECEPTOR
 (ERYTHROPOIETIN(W)RECEPTOR)
 23 (EPO OR (ERYTHROPOIETIN RECEPTOR))
 FILE 'SCISEARCH'

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4531 EPO
222 EPOS
4734 EPO
      (EPO OR EPOS)
16279 ERYTHROPOIETIN
46 ERYTHROPOIETINS
16295 ERYTHROPOIETIN
      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
583687 RECEPTOR
306298 RECEPTORS
721251 RECEPTOR
      (RECEPTOR OR RECEPTORS)
1592 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
5851 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'SYNTHLINE'
1 EPO
0 "ERYTHROPOIETIN"
798 "RECEPTOR"
103 "RECEPTORS"
876 "RECEPTOR"
      ("RECEPTOR" OR "RECEPTORS")
0 ERYTHROPOIETIN RECEPTOR
      ("ERYTHROPOIETIN(W)RECEPTOR")
1 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'TOXCENTER'
2126 EPO
34 EPOS
2145 EPO
      (EPO OR EPOS)
6597 ERYTHROPOIETIN
72 ERYTHROPOIETINS
6605 ERYTHROPOIETIN
      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
264727 RECEPTOR
166955 RECEPTORS
322116 RECEPTOR
      (RECEPTOR OR RECEPTORS)
285 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
2274 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'USPATFULL'
13847 EPO
184 EPOS
13982 EPO
      (EPO OR EPOS)
7989 ERYTHROPOIETIN
1884 ERYTHROPOIETINS
9673 ERYTHROPOIETIN
      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
104220 RECEPTOR
75854 RECEPTORS
118925 RECEPTOR
      (RECEPTOR OR RECEPTORS)
773 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
14435 (EPO OR (ERYTHROPOIETIN RECEPTOR))

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      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
41921 RECEPTOR
13906 RECEPTORS
46849 RECEPTOR
      (RECEPTOR OR RECEPTORS)
80 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
659 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'WPIFV'
10 EPO
15 ERYTHROPOIETIN
354 RECEPTOR
89 RECEPTORS
389 RECEPTOR
      (RECEPTOR OR RECEPTORS)
2 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
10 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'WPINDEX'
580 EPO
36 EPOS
612 EPO
      (EPO OR EPOS)
1446 ERYTHROPOIETIN
16 ERYTHROPOIETINS
1458 ERYTHROPOIETIN
      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
41921 RECEPTOR
13906 RECEPTORS
46849 RECEPTOR
      (RECEPTOR OR RECEPTORS)
80 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
659 (EPO OR (ERYTHROPOIETIN RECEPTOR))

```

L8 QUE (EPO OR (ERYTHROPOIETIN RECEPTOR))

=> DIS HIST

(FILE 'HOME' ENTERED AT 10:25:10 ON 28 DEC 2004)

FILE 'REGISTRY' ENTERED AT 10:25:22 ON 28 DEC 2004

```

L1 STRUCTURE UPLOADED
L2 0 S L1 SAM
L3 0 S L1 FAM
L4 8 S L1 FUL
L5 1 S L1 FUL FAM
L6 0 L5 NOT L4

```

FILE 'HCAPLUS, USPATFULL, CHEMCATS' ENTERED AT 10:35:54 ON 28 DEC 2004

INDEX 'ADISCTI, ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, ANTE, AQUALINE, AQUASCI, BIOBUSINESS, BIOCOMMERCE, BIOENG, BIOSIS, BIOTECHABS, BIOTECHDS,

```

FILE 'USPAT2'
807 EPO
13 EPOS
818 EPO
      (EPO OR EPOS)
477 ERYTHROPOIETIN
30 ERYTHROPOIETINS
494 ERYTHROPOIETIN
      (ERYTHROPOIETIN OR ERYTHROPOIETINS)
7293 RECEPTOR
5119 RECEPTORS
8267 RECEPTOR
      (RECEPTOR OR RECEPTORS)
32 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
836 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'VETB'
0 EPO
1 ERYTHROPOIETIN
77 RECEPTOR
25 RECEPTORS
82 RECEPTOR
      (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
0 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'VETU'
23 EPO
65 ERYTHROPOIETIN
970 RECEPTOR
641 RECEPTORS
1302 RECEPTOR
      (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
23 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'WATER'
0 EPO
2 EPOS
2 EPO
      (EPO OR EPOS)
1 ERYTHROPOIETIN
471 RECEPTOR
265 RECEPTORS
641 RECEPTOR
      (RECEPTOR OR RECEPTORS)
0 ERYTHROPOIETIN RECEPTOR
      (ERYTHROPOIETIN(W)RECEPTOR)
2 (EPO OR (ERYTHROPOIETIN RECEPTOR))
FILE 'WPIDS'
580 EPO
36 EPOS
612 EPO
      (EPO OR EPOS)
1446 ERYTHROPOIETIN
16 ERYTHROPOIETINS
1458 ERYTHROPOIETIN

```

BIOTECHNO, CABA, CANCERLIT, CAPLUS, CEABA-VTB, CEN, CIN, CONFSCI, CROPB, CROPU, DDFB, DDFU, DGENE, DISSABS, ...' ENTERED AT 10:38:10 ON 28 DEC 2004

SEA (EPO OR (ERYTHROPOIETIN RECEPTOR))

```

73 FILE ADISCTI
27 FILE ADISINSIGHT
19 FILE ADISNEWS
63 FILE AGRICOLA
53 FILE ANABSTR
22 FILE ANTE
3 FILE AQUALINE
99 FILE AQUASCI
410 FILE BIOBUSINESS
1197 FILE BIOCOMMERCE
199 FILE BIOENG
6546 FILE BIOSIS
407 FILE BIOTECHABS
407 FILE BIOTECHDS
2960 FILE BIOTECHNO
308 FILE CABA
3775 FILE CANCERLIT
6216 FILE CAPLUS
124 FILE CEABA-VTB
34 FILE CEN
539 FILE CIN
165 FILE CONFSCI
9 FILE CROPB
67 FILE CROPU
3 FILE DDFB
1441 FILE DDFU
15707 FILE DGENE
216 FILE DISSABS
3 FILE DRUGB
76 FILE DRUGMONOG2
1657 FILE DRUGU
57 FILE EMBAL
5095 FILE EMBASE
2150 FILE ESBIODBASE
139 FILE FEDRIP
13 FILE FOMAD
29 FILE FROSTI
25 FILE FSTA
6457 FILE GENBANK
7 FILE HEALSAFE
74 FILE IFIPAT
74 FILE IMSDRUGNEWS
33 FILE IMSPRODUCT
36 FILE IMSRESEARCH
800 FILE JICST-EPLUS
6 FILE KOSMET
1047 FILE LIFESCI
11 FILE MEDICONF
5281 FILE MEDLINE
15 FILE NIOSHTIC
108 FILE NTIS

```

```

11 FILE NUTRACEUT
75 FILE OCEAN
2342 FILE PASCAL
73 FILE PHAR
101 FILE PHARMAML
7 FILE PHIC
861 FILE PHIN
4612 FILE PROMT
22 FILE PROUSDDR
1 FILE PS
23 FILE RDISCLOSURE
5851 FILE SCISEARCH
1 FILE SYNTHLINE
2274 FILE TOXCENTER
14435 FILE USPATFULL
836 FILE USPAT2
23 FILE VETU
2 FILE WATER
659 FILE WPIDS
10 FILE WPIFV
659 FILE WPINDE
L8 QUE (EPO OR (ERYTHROPOIETIN RECEPTOR))
-----

```

=>

---Logging off of STN---

=>

Executing the logoff script...

=> LOG Y

| | |
|--|------------|
| COST IN U.S. DOLLARS | SINCE FILE |
| TOTAL | ENTRY |
| SESSION | |
| FULL ESTIMATED COST | 11.97 |
| 325.86 | |
| DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) | SINCE FILE |
| TOTAL | ENTRY |
| SESSION | |
| CA SUBSCRIBER PRICE | 0.00 |
| 1.32 | |

STN INTERNATIONAL LOGOFF AT 10:50:37 ON 28 DEC 2004